



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 04:29 PM GMT

PDB ID : 4F2M
Title : Crystal structure of a TGEV coronavirus Spike fragment in complex with the TGEV neutralizing monoclonal antibody 1AF10
Authors : Reguera, J.; Santiago, C.; Mudgal, G.; Ordone, D.; Enjuanes, L.; Casasnovas, J.M.
Deposited on : 2012-05-08
Resolution : 3.00 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

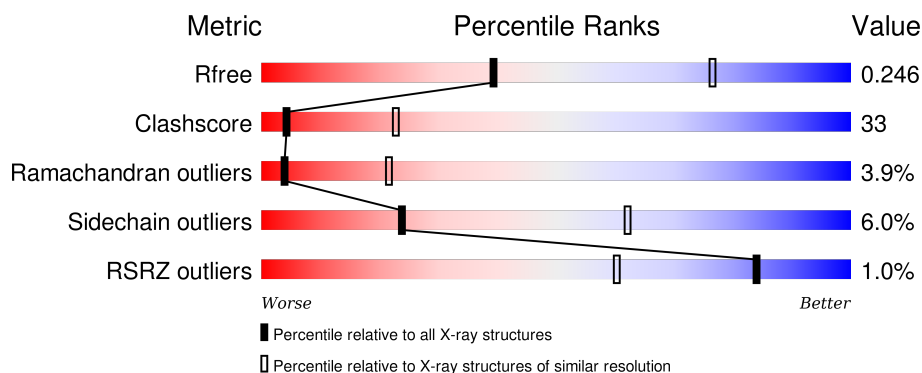
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	1578 (3.00-3.00)
Clashscore	102246	1912 (3.00-3.00)
Ramachandran outliers	100387	1853 (3.00-3.00)
Sidechain outliers	100360	1856 (3.00-3.00)
RSRZ outliers	91569	1592 (3.00-3.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	221	<div> <div></div> <div>51% 41% 5% .</div> </div>
1	C	221	<div> <div>%</div> <div>47% 45% 6% .</div> </div>
2	B	214	<div> <div>%</div> <div>51% 41% 6% .</div> </div>
2	D	214	<div> <div></div> <div>57% 39% ..</div> </div>
3	E	209	<div> <div>%</div> <div>33% 32% 5% . 29%</div> </div>

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Mol	Chain	Length	Quality of chain
3	F	209	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	ACY	B	301	-	-	-	X
5	NAG	F	701	-	-	-	X
6	NAG	F	705	X	-	-	-

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 8998 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called monoclonal antibody 1AF10, heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	214	Total	C	N	O	S	0	0	0
			1614	1024	264	319	7			
1	C	220	Total	C	N	O	S	1	0	0
			1655	1047	270	330	8			

- Molecule 2 is a protein called monoclonal antibody 1AF10, light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	213	Total	C	N	O	S	1	0	0
			1643	1017	280	341	5			
2	D	211	Total	C	N	O	S	0	0	0
			1626	1008	277	336	5			

- Molecule 3 is a protein called Spike protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	148	Total	C	N	O	S	0	0	0
			1162	736	194	224	8			
3	F	148	Total	C	N	O	S	0	0	0
			1162	736	194	224	8			

There are 78 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	456	TYR	-	EXPRESSION TAG	UNP Q0PKZ5
E	457	PRO	-	EXPRESSION TAG	UNP Q0PKZ5
E	458	TYR	-	EXPRESSION TAG	UNP Q0PKZ5
E	459	ASP	-	EXPRESSION TAG	UNP Q0PKZ5
E	460	VAL	-	EXPRESSION TAG	UNP Q0PKZ5
E	461	PRO	-	EXPRESSION TAG	UNP Q0PKZ5
E	462	ASP	-	EXPRESSION TAG	UNP Q0PKZ5
E	463	TYR	-	EXPRESSION TAG	UNP Q0PKZ5

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Chain	Residue	Modelled	Actual	Comment	Reference
E	464	ALA	-	EXPRESSION TAG	UNP Q0PKZ5
E	465	GLY	-	EXPRESSION TAG	UNP Q0PKZ5
E	466	ALA	-	EXPRESSION TAG	UNP Q0PKZ5
E	467	GLN	-	EXPRESSION TAG	UNP Q0PKZ5
E	468	PRO	-	EXPRESSION TAG	UNP Q0PKZ5
E	469	ALA	-	EXPRESSION TAG	UNP Q0PKZ5
E	470	ARG	-	EXPRESSION TAG	UNP Q0PKZ5
E	471	SER	-	EXPRESSION TAG	UNP Q0PKZ5
E	472	PRO	-	EXPRESSION TAG	UNP Q0PKZ5
E	473	GLY	-	EXPRESSION TAG	UNP Q0PKZ5
E	474	LEU	-	EXPRESSION TAG	UNP Q0PKZ5
E	475	VAL	-	EXPRESSION TAG	UNP Q0PKZ5
E	476	PRO	-	EXPRESSION TAG	UNP Q0PKZ5
E	477	ARG	-	EXPRESSION TAG	UNP Q0PKZ5
E	478	GLY	-	EXPRESSION TAG	UNP Q0PKZ5
E	479	SER	-	EXPRESSION TAG	UNP Q0PKZ5
E	480	ARG	-	EXPRESSION TAG	UNP Q0PKZ5
E	651	LEU	-	EXPRESSION TAG	UNP Q0PKZ5
E	652	VAL	-	EXPRESSION TAG	UNP Q0PKZ5
E	653	PRO	-	EXPRESSION TAG	UNP Q0PKZ5
E	654	ARG	-	EXPRESSION TAG	UNP Q0PKZ5
E	655	GLY	-	EXPRESSION TAG	UNP Q0PKZ5
E	656	SER	-	EXPRESSION TAG	UNP Q0PKZ5
E	657	ASP	-	EXPRESSION TAG	UNP Q0PKZ5
E	658	TYR	-	EXPRESSION TAG	UNP Q0PKZ5
E	659	LYS	-	EXPRESSION TAG	UNP Q0PKZ5
E	660	ASP	-	EXPRESSION TAG	UNP Q0PKZ5
E	661	ASP	-	EXPRESSION TAG	UNP Q0PKZ5
E	662	ASP	-	EXPRESSION TAG	UNP Q0PKZ5
E	663	ASP	-	EXPRESSION TAG	UNP Q0PKZ5
E	664	LYS	-	EXPRESSION TAG	UNP Q0PKZ5
F	456	TYR	-	EXPRESSION TAG	UNP Q0PKZ5
F	457	PRO	-	EXPRESSION TAG	UNP Q0PKZ5
F	458	TYR	-	EXPRESSION TAG	UNP Q0PKZ5
F	459	ASP	-	EXPRESSION TAG	UNP Q0PKZ5
F	460	VAL	-	EXPRESSION TAG	UNP Q0PKZ5
F	461	PRO	-	EXPRESSION TAG	UNP Q0PKZ5
F	462	ASP	-	EXPRESSION TAG	UNP Q0PKZ5
F	463	TYR	-	EXPRESSION TAG	UNP Q0PKZ5
F	464	ALA	-	EXPRESSION TAG	UNP Q0PKZ5
F	465	GLY	-	EXPRESSION TAG	UNP Q0PKZ5
F	466	ALA	-	EXPRESSION TAG	UNP Q0PKZ5

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Chain	Residue	Modelled	Actual	Comment	Reference
F	467	GLN	-	EXPRESSION TAG	UNP Q0PKZ5
F	468	PRO	-	EXPRESSION TAG	UNP Q0PKZ5
F	469	ALA	-	EXPRESSION TAG	UNP Q0PKZ5
F	470	ARG	-	EXPRESSION TAG	UNP Q0PKZ5
F	471	SER	-	EXPRESSION TAG	UNP Q0PKZ5
F	472	PRO	-	EXPRESSION TAG	UNP Q0PKZ5
F	473	GLY	-	EXPRESSION TAG	UNP Q0PKZ5
F	474	LEU	-	EXPRESSION TAG	UNP Q0PKZ5
F	475	VAL	-	EXPRESSION TAG	UNP Q0PKZ5
F	476	PRO	-	EXPRESSION TAG	UNP Q0PKZ5
F	477	ARG	-	EXPRESSION TAG	UNP Q0PKZ5
F	478	GLY	-	EXPRESSION TAG	UNP Q0PKZ5
F	479	SER	-	EXPRESSION TAG	UNP Q0PKZ5
F	480	ARG	-	EXPRESSION TAG	UNP Q0PKZ5
F	651	LEU	-	EXPRESSION TAG	UNP Q0PKZ5
F	652	VAL	-	EXPRESSION TAG	UNP Q0PKZ5
F	653	PRO	-	EXPRESSION TAG	UNP Q0PKZ5
F	654	ARG	-	EXPRESSION TAG	UNP Q0PKZ5
F	655	GLY	-	EXPRESSION TAG	UNP Q0PKZ5
F	656	SER	-	EXPRESSION TAG	UNP Q0PKZ5
F	657	ASP	-	EXPRESSION TAG	UNP Q0PKZ5
F	658	TYR	-	EXPRESSION TAG	UNP Q0PKZ5
F	659	LYS	-	EXPRESSION TAG	UNP Q0PKZ5
F	660	ASP	-	EXPRESSION TAG	UNP Q0PKZ5
F	661	ASP	-	EXPRESSION TAG	UNP Q0PKZ5
F	662	ASP	-	EXPRESSION TAG	UNP Q0PKZ5
F	663	ASP	-	EXPRESSION TAG	UNP Q0PKZ5
F	664	LYS	-	EXPRESSION TAG	UNP Q0PKZ5

- Molecule 4 is ACETIC ACID (three-letter code: ACY) (formula: C₂H₄O₂).

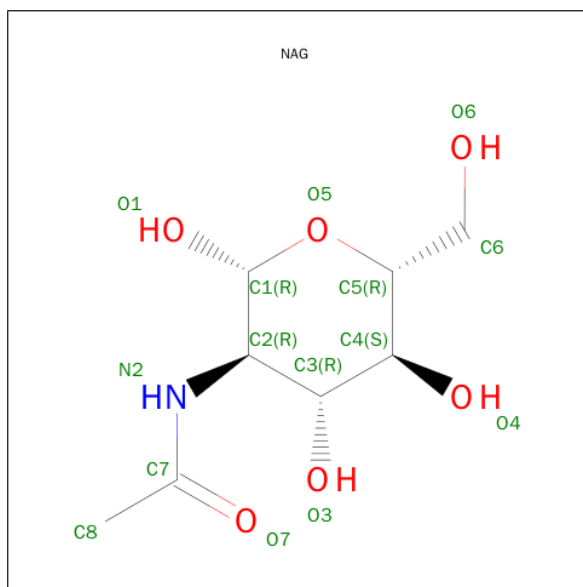


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			4	2	2		
4	D	1	Total	C	O	0	0
			4	2	2		

- Molecule 5 is a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	E	2	Total	C	N	O	0	0
			28	16	2	10		
5	E	2	Total	C	N	O	0	0
			28	16	2	10		
5	F	2	Total	C	N	O	0	0
			28	16	2	10		
5	F	2	Total	C	N	O	0	0
			28	16	2	10		

- Molecule 6 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	F	1	Total	C	N	O	0	0
			14	8	1	5		

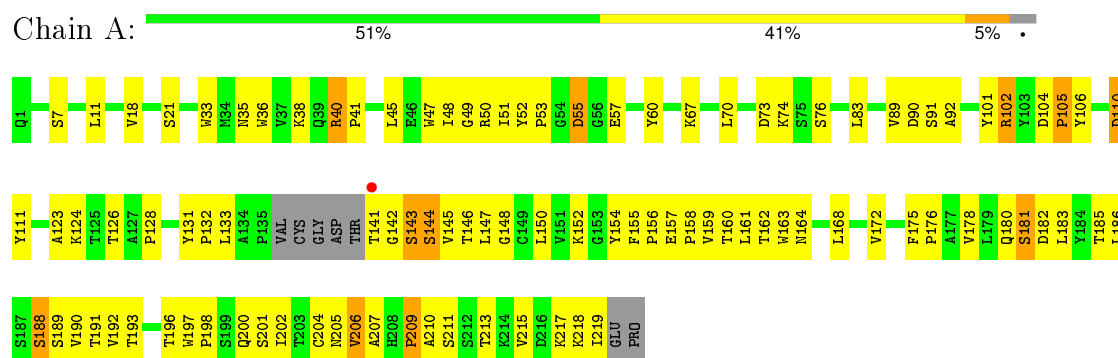
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	B	1	Total	O	0	0
			1	1		
7	D	1	Total	O	0	0
			1	1		

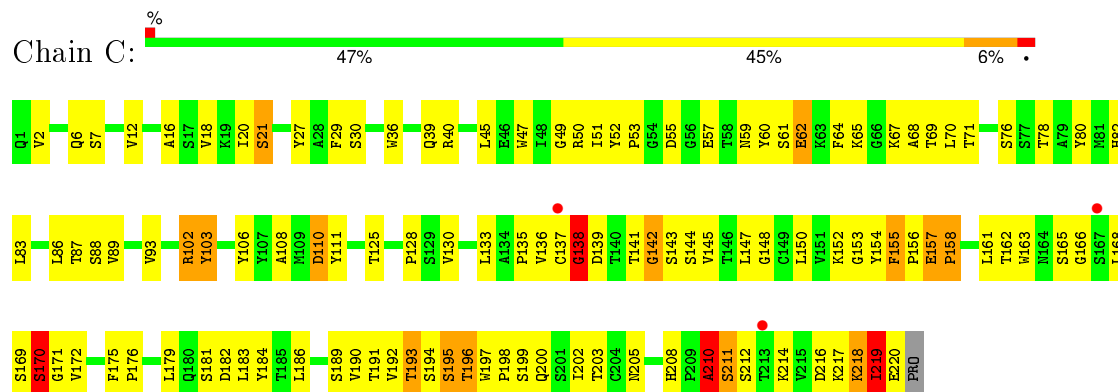
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($\text{RSRZ} > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

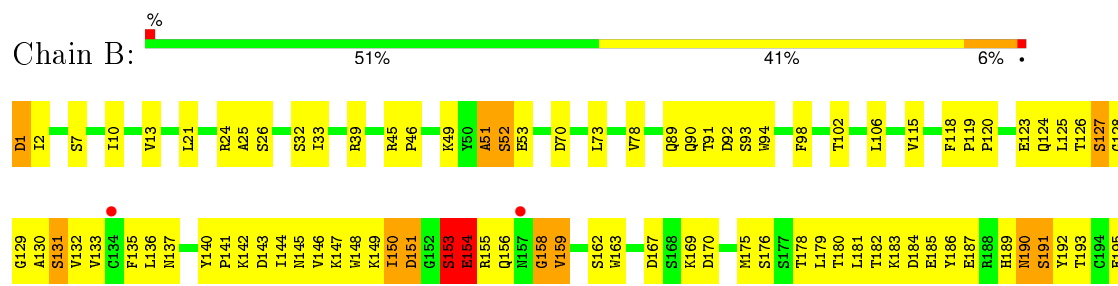
- Molecule 1: monoclonal antibody 1AF10, heavy chain



- Molecule 1: monoclonal antibody 1AF10, heavy chain



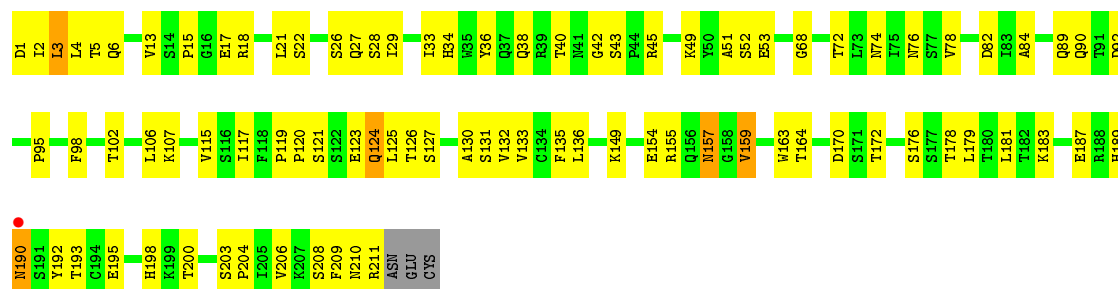
- Molecule 2: monoclonal antibody 1AF10, light chain





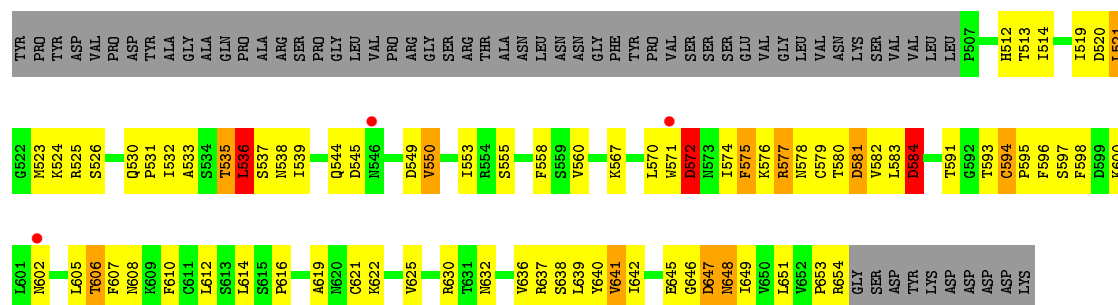
- Molecule 2: monoclonal antibody 1AF10, light chain

Chain D: 57% 39% ..



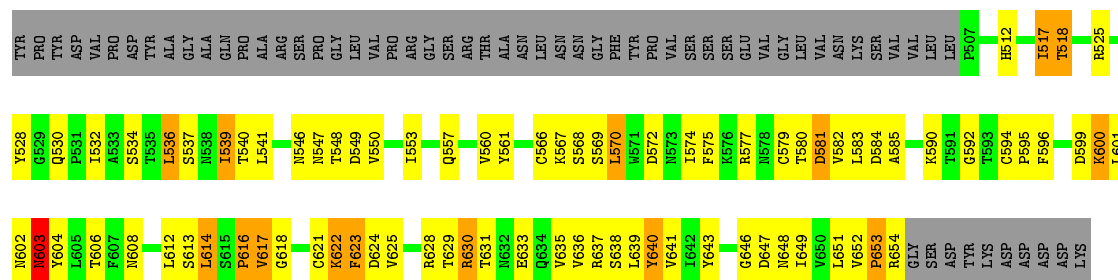
- Molecule 3: Spike protein

Chain E: 33% 32% 5% 29%



- Molecule 3: Spike protein

Chain F: 31% 32% 7% 29%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	82.77Å 106.05Å 103.52Å 90.00° 95.13° 90.00°	Depositor
Resolution (Å)	25.00 – 3.00 24.64 – 3.00	Depositor EDS
% Data completeness (in resolution range)	94.7 (25.00-3.00) 94.8 (24.64-3.00)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.11 (at 2.99Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.4_486)	Depositor
R, R_{free}	0.216 , 0.251 0.209 , 0.246	Depositor DCC
R_{free} test set	3095 reflections (5.05%)	DCC
Wilson B-factor (Å ²)	67.6	Xtriage
Anisotropy	0.324	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 47.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 33910 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	8998	wwPDB-VP
Average B, all atoms (Å ²)	82.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.31% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ACY, NAG

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.28	0/1658	0.52	0/2261
1	C	0.32	0/1700	0.63	1/2320 (0.0%)
2	B	0.33	0/1679	0.63	1/2282 (0.0%)
2	D	0.29	0/1662	0.49	0/2259
3	E	0.32	0/1186	0.58	0/1614
3	F	0.36	0/1186	0.63	0/1614
All	All	0.31	0/9071	0.58	2/12350 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	C	0	4
2	B	0	3
3	E	0	3
3	F	0	2
All	All	0	12

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	138	GLY	N-CA-C	-6.79	96.12	113.10
2	B	212	ASN	N-CA-C	5.45	125.72	111.00

There are no chirality outliers.

All (12) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	150	ILE	Peptide
2	B	158	GLY	Peptide
2	B	211	ARG	Peptide
1	C	138	GLY	Peptide
1	C	157	GLU	Peptide
1	C	166	GLY	Peptide
1	C	210	ALA	Peptide
3	E	536	LEU	Peptide
3	E	572	ASP	Peptide
3	E	600	LYS	Peptide
3	F	616	PRO	Peptide
3	F	622	LYS	Peptide

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1614	0	1568	97	0
1	C	1655	0	1603	130	0
2	B	1643	0	1570	120	0
2	D	1626	0	1558	67	0
3	E	1162	0	1133	79	0
3	F	1162	0	1132	120	0
4	B	4	0	3	1	0
4	D	4	0	3	0	0
5	E	56	0	50	0	0
5	F	56	0	50	3	0
6	F	14	0	13	0	0
7	B	1	0	0	0	0
7	D	1	0	0	1	0
All	All	8998	0	8683	586	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 33.

All (586) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:521:LEU:HA	3:E:537:SER:HB3	1.22	1.18
3:F:590:LYS:O	3:F:623:PHE:HB2	1.52	1.09
3:F:616:PRO:HG2	3:F:640:TYR:CD2	1.89	1.06
3:F:549:ASP:HA	3:F:614:LEU:HB2	1.37	1.02
2:B:151:ASP:CG	2:B:190:ASN:HB2	1.82	0.99
2:B:187:GLU:HA	2:B:211:ARG:HH22	1.25	0.99
3:F:630:ARG:N	3:F:630:ARG:HE	1.62	0.98
2:B:195:GLU:HG2	2:B:206:VAL:HG12	1.46	0.97
1:C:30:SER:HA	1:C:53:PRO:HB2	1.47	0.96
2:B:150:ILE:HG22	2:B:190:ASN:OD1	1.65	0.94
1:C:40:ARG:NH2	1:C:89:VAL:HG13	1.83	0.94
3:F:614:LEU:O	3:F:616:PRO:HD3	1.68	0.92
1:A:142:GLY:N	1:A:143:SER:HA	1.84	0.91
2:B:189:HIS:HE1	2:B:210:ASN:HA	1.34	0.91
2:B:189:HIS:NE2	2:B:210:ASN:HB2	1.86	0.90
2:B:118:PHE:HB2	2:B:133:VAL:HG22	1.54	0.90
2:B:144:ILE:HG12	2:B:145:ASN:H	1.38	0.88
2:B:189:HIS:CE1	2:B:210:ASN:HA	2.10	0.87
1:C:155:PHE:CD1	1:C:156:PRO:HA	2.09	0.87
1:C:50:ARG:HE	1:C:59:ASN:HD22	1.19	0.87
3:E:519:ILE:HG23	3:E:539:ILE:HG22	1.57	0.86
2:D:193:THR:HG22	2:D:208:SER:HB3	1.58	0.84
3:F:580:THR:HG22	3:F:581:ASP:H	1.42	0.83
1:C:136:VAL:HG11	2:D:119:PRO:HG2	1.59	0.83
2:D:21:LEU:HD22	2:D:102:THR:HG21	1.60	0.83
1:C:125:THR:HA	1:C:156:PRO:HD3	1.60	0.83
3:F:616:PRO:HG2	3:F:640:TYR:CG	2.12	0.83
3:F:592:GLY:H	3:F:623:PHE:HB3	1.45	0.81
1:A:141:THR:HG23	1:A:143:SER:HB2	1.61	0.81
2:D:190:ASN:HB2	2:D:211:ARG:HD3	1.62	0.81
3:F:640:TYR:HD2	3:F:640:TYR:H	1.26	0.80
2:B:198:HIS:CD2	2:B:200:THR:HB	2.16	0.80
3:E:521:LEU:HA	3:E:537:SER:CB	2.08	0.80
3:F:624:ASP:HB3	3:F:635:VAL:HG13	1.62	0.79
1:C:50:ARG:NE	1:C:59:ASN:HD22	1.81	0.78
1:C:218:LYS:O	1:C:219:ILE:HG23	1.84	0.78
2:B:180:THR:O	2:B:181:LEU:HD12	1.82	0.78
1:C:125:THR:HA	1:C:156:PRO:CD	2.14	0.78
3:F:536:LEU:HD13	3:F:536:LEU:H	1.49	0.77
3:F:539:ILE:HG23	3:F:614:LEU:HD11	1.66	0.77
2:D:33:ILE:HG22	2:D:51:ALA:HB2	1.67	0.77
3:F:570:LEU:HD21	3:F:577:ARG:HH21	1.51	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:567:LYS:HB3	3:E:575:PHE:O	1.85	0.76
2:B:150:ILE:HA	2:B:191:SER:O	1.86	0.75
1:C:55:ASP:OD2	1:C:57:GLU:HG2	1.87	0.74
3:F:628:ARG:HH11	3:F:628:ARG:HG2	1.52	0.74
1:A:40:ARG:NH1	1:A:91:SER:O	2.21	0.74
2:B:147:LYS:HZ1	2:B:154:GLU:HB3	1.53	0.73
2:D:190:ASN:OD1	2:D:211:ARG:N	2.21	0.72
1:C:125:THR:HG23	1:C:156:PRO:HD2	1.71	0.72
1:A:7:SER:HB3	1:A:21:SER:HB3	1.71	0.72
3:F:640:TYR:N	3:F:640:TYR:HD2	1.87	0.72
1:A:40:ARG:HG3	1:A:41:PRO:HD2	1.72	0.72
2:B:2:ILE:HD11	2:B:25:ALA:HB1	1.71	0.72
3:F:640:TYR:N	3:F:640:TYR:CD2	2.58	0.72
2:B:148:TRP:CD2	2:B:179:LEU:HD12	2.25	0.71
3:F:612:LEU:HD22	3:F:639:LEU:HD11	1.73	0.71
2:B:153:SER:HB2	2:B:154:GLU:OE2	1.90	0.71
1:C:29:PHE:O	1:C:53:PRO:HG2	1.91	0.70
2:B:13:VAL:HG21	2:B:78:VAL:HG21	1.74	0.70
1:A:60:TYR:CE1	1:A:70:LEU:HD13	2.26	0.70
1:C:163:TRP:HZ3	1:C:218:LYS:NZ	1.89	0.70
1:A:172:VAL:HA	1:A:189:SER:O	1.91	0.70
2:B:193:THR:HG22	2:B:208:SER:HB3	1.74	0.69
1:C:40:ARG:HH22	1:C:89:VAL:HG13	1.55	0.69
3:F:512:HIS:ND1	3:F:557:GLN:HG2	2.08	0.69
1:A:217:LYS:HG2	1:A:218:LYS:H	1.58	0.69
3:E:574:ILE:HG23	3:E:574:ILE:O	1.92	0.69
2:B:179:LEU:HD23	2:B:181:LEU:HD13	1.74	0.69
1:A:209:PRO:C	1:A:211:SER:H	1.97	0.68
1:C:203:THR:HA	1:C:216:ASP:O	1.93	0.68
3:E:560:VAL:HG21	3:E:598:PHE:CZ	2.28	0.68
3:E:653:PRO:HA	3:E:654:ARG:CB	2.22	0.68
2:D:195:GLU:HG3	2:D:206:VAL:HG12	1.74	0.68
2:B:118:PHE:HB2	2:B:133:VAL:CG2	2.23	0.67
3:F:603:ASN:O	3:F:604:TYR:CG	2.47	0.67
1:A:157:GLU:HG2	1:A:158:PRO:HA	1.76	0.67
3:E:610:PHE:CE2	3:E:612:LEU:HD11	2.29	0.67
1:C:197:TRP:CD1	1:C:202:ILE:HD12	2.29	0.67
1:A:202:ILE:HG13	1:A:218:LYS:HE3	1.76	0.67
3:F:599:ASP:O	3:F:600:LYS:HB2	1.94	0.67
3:F:614:LEU:C	3:F:616:PRO:HD3	2.14	0.66
2:B:147:LYS:NZ	2:B:154:GLU:HB3	2.09	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:190:ASN:HA	2:D:192:TYR:CE1	2.31	0.66
1:C:157:GLU:HG2	1:C:158:PRO:HA	1.77	0.66
1:A:55:ASP:N	1:A:55:ASP:OD1	2.29	0.66
2:B:49:LYS:O	2:B:53:GLU:HB3	1.95	0.66
2:B:132:VAL:CG1	2:B:179:LEU:HB3	2.26	0.66
1:C:203:THR:HG22	1:C:217:LYS:HG2	1.79	0.65
1:A:196:THR:O	1:A:200:GLN:N	2.27	0.65
1:A:192:VAL:HG22	1:A:193:THR:H	1.60	0.65
3:F:548:THR:HG23	3:F:549:ASP:OD1	1.96	0.65
2:B:189:HIS:HB2	2:B:211:ARG:HD2	1.78	0.65
1:C:219:ILE:HD12	1:C:220:GLU:H	1.62	0.65
3:F:548:THR:O	3:F:549:ASP:HB2	1.97	0.65
2:B:133:VAL:HG12	2:B:178:THR:HG23	1.79	0.65
1:C:102:ARG:O	1:C:103:TYR:HB2	1.97	0.65
1:C:219:ILE:O	1:C:220:GLU:HB2	1.96	0.64
2:D:159:VAL:HG12	2:D:179:LEU:HD13	1.80	0.64
3:F:590:LYS:O	3:F:623:PHE:CB	2.40	0.64
1:C:50:ARG:HE	1:C:59:ASN:ND2	1.92	0.64
2:D:49:LYS:O	2:D:53:GLU:HB3	1.97	0.64
1:C:12:VAL:HG11	1:C:18:VAL:HG12	1.81	0.63
1:C:40:ARG:HH21	1:C:89:VAL:HG13	1.62	0.63
1:C:219:ILE:HG13	1:C:220:GLU:N	2.13	0.63
3:F:622:LYS:HE2	3:F:638:SER:CB	2.29	0.63
3:F:638:SER:HB2	3:F:640:TYR:CE2	2.33	0.63
2:B:132:VAL:HG13	2:B:179:LEU:HB3	1.79	0.63
2:B:181:LEU:HG	2:B:185:GLU:OE2	1.97	0.63
2:B:24:ARG:HG2	2:B:70:ASP:OD1	1.99	0.63
2:B:167:ASP:OD2	2:B:169:LYS:HB3	1.98	0.63
2:B:33:ILE:HG23	2:B:51:ALA:HB2	1.81	0.63
1:C:52:TYR:CE2	1:C:102:ARG:HB2	2.33	0.62
2:B:90:GLN:HE21	2:B:92:ASP:N	1.96	0.62
2:B:90:GLN:NE2	2:B:93:SER:H	1.98	0.62
2:B:198:HIS:HD2	2:B:200:THR:HB	1.63	0.62
3:E:514:ILE:CG2	3:F:512:HIS:HB3	2.30	0.62
3:F:537:SER:HB2	3:F:636:VAL:HG21	1.82	0.62
2:B:179:LEU:HD23	2:B:181:LEU:CD1	2.28	0.62
3:E:560:VAL:HG21	3:E:598:PHE:HZ	1.65	0.62
3:E:535:THR:O	3:E:536:LEU:HB2	2.00	0.62
2:B:115:VAL:HG12	2:B:136:LEU:HG	1.81	0.61
3:F:541:LEU:HD13	3:F:614:LEU:HD21	1.81	0.61
1:C:172:VAL:HA	1:C:189:SER:O	2.00	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:622:LYS:HA	3:F:640:TYR:HA	1.82	0.61
1:A:60:TYR:HE1	1:A:70:LEU:HD13	1.65	0.61
1:A:197:TRP:CD1	1:A:202:ILE:HG12	2.36	0.61
2:D:126:THR:HG22	2:D:126:THR:O	2.00	0.61
1:C:157:GLU:HG2	1:C:158:PRO:CA	2.31	0.61
5:F:701:NAG:H62	5:F:702:NAG:C7	2.31	0.61
1:C:102:ARG:HG3	1:C:102:ARG:HH11	1.66	0.60
1:C:176:PRO:HD3	2:D:164:THR:HG22	1.83	0.60
3:F:568:SER:HB3	3:F:577:ARG:CB	2.31	0.60
3:F:572:ASP:OD2	3:F:574:ILE:HG22	2.02	0.60
1:C:192:VAL:O	1:C:192:VAL:HG13	2.02	0.60
1:C:163:TRP:CZ3	1:C:218:LYS:NZ	2.69	0.60
1:C:71:THR:OG1	1:C:80:TYR:HB2	2.02	0.60
2:B:144:ILE:HG12	2:B:145:ASN:N	2.14	0.59
1:A:160:THR:CG2	1:A:207:ALA:HB3	2.33	0.59
1:A:141:THR:CG2	1:A:143:SER:HB2	2.31	0.59
2:B:181:LEU:HD23	2:B:185:GLU:HG2	1.83	0.59
1:A:144:SER:HB3	1:A:191:THR:HG22	1.85	0.59
1:A:145:VAL:N	1:A:192:VAL:O	2.35	0.59
1:A:123:ALA:HB2	1:A:182:ASP:OD1	2.01	0.59
1:A:161:LEU:HD13	1:A:206:VAL:HG13	1.83	0.59
2:D:115:VAL:HG12	2:D:136:LEU:HG	1.85	0.59
3:E:579:CYS:O	3:E:584:ASP:HA	2.02	0.59
1:C:141:THR:O	1:C:142:GLY:C	2.40	0.59
2:B:212:ASN:O	2:B:213:GLU:HG3	2.03	0.58
3:E:610:PHE:HE2	3:E:612:LEU:HD11	1.68	0.58
2:B:10:ILE:HD13	2:B:142:LYS:NZ	2.18	0.58
3:F:580:THR:HG22	3:F:581:ASP:N	2.14	0.58
1:C:136:VAL:C	1:C:138:GLY:HA2	2.23	0.58
1:C:208:HIS:CE1	1:C:210:ALA:HB3	2.39	0.58
1:C:203:THR:HG22	1:C:217:LYS:CG	2.34	0.58
2:B:195:GLU:HG2	2:B:206:VAL:CG1	2.30	0.58
1:C:182:ASP:O	1:C:183:LEU:HD13	2.04	0.58
2:B:127:SER:C	2:B:129:GLY:N	2.56	0.58
2:D:28:SER:HA	2:D:68:GLY:O	2.04	0.58
3:E:581:ASP:OD2	3:E:582:VAL:HG22	2.02	0.58
2:D:119:PRO:HG3	2:D:209:PHE:CD2	2.39	0.57
2:B:90:GLN:HE21	2:B:92:ASP:H	1.52	0.57
1:C:219:ILE:CG1	1:C:220:GLU:N	2.67	0.57
3:E:572:ASP:HB2	3:E:574:ILE:HG22	1.86	0.57
1:A:192:VAL:HG22	1:A:193:THR:N	2.19	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:550:VAL:HA	3:F:612:LEU:O	2.04	0.57
1:C:218:LYS:C	1:C:219:ILE:HG23	2.23	0.57
2:B:136:LEU:HD22	2:B:175:MET:CE	2.34	0.57
3:F:625:VAL:HG23	3:F:636:VAL:HG13	1.86	0.57
2:B:183:LYS:O	2:B:187:GLU:N	2.37	0.57
3:F:568:SER:HB3	3:F:577:ARG:HB2	1.87	0.57
1:C:136:VAL:O	1:C:138:GLY:N	2.38	0.57
2:B:148:TRP:CG	2:B:179:LEU:HD12	2.39	0.56
3:E:524:LYS:HG2	3:E:525:ARG:N	2.20	0.56
3:E:612:LEU:HD12	3:E:612:LEU:N	2.20	0.56
3:F:561:TYR:CE2	5:F:701:NAG:H82	2.39	0.56
2:B:125:LEU:HA	2:B:129:GLY:O	2.05	0.56
2:B:189:HIS:CE1	2:B:210:ASN:CA	2.85	0.56
1:C:163:TRP:HZ3	1:C:218:LYS:HZ2	1.51	0.56
1:A:133:LEU:HB2	1:A:148:GLY:C	2.26	0.56
1:A:106:TYR:OH	3:E:532:ILE:HD13	2.05	0.56
2:B:148:TRP:CD1	2:B:159:VAL:HG11	2.41	0.56
2:B:182:THR:OG1	2:B:185:GLU:HB3	2.05	0.56
1:C:16:ALA:O	1:C:86:LEU:HD13	2.05	0.56
2:B:189:HIS:CG	2:B:211:ARG:HD2	2.40	0.56
2:D:89:GLN:HB2	2:D:98:PHE:CE1	2.41	0.56
3:F:608:ASN:CG	3:F:646:GLY:HA3	2.26	0.56
3:F:629:THR:C	3:F:630:ARG:HE	2.07	0.56
2:D:121:SER:HB2	2:D:123:GLU:OE1	2.06	0.56
3:F:541:LEU:HD13	3:F:614:LEU:CD2	2.35	0.56
1:A:133:LEU:HD11	1:A:150:LEU:HB2	1.88	0.56
2:B:21:LEU:HD22	2:B:102:THR:HG21	1.88	0.56
1:A:186:LEU:HD12	1:A:186:LEU:C	2.26	0.55
1:C:62:GLU:HA	1:C:65:LYS:HG3	1.88	0.55
1:C:155:PHE:CG	1:C:156:PRO:HA	2.41	0.55
3:F:592:GLY:HA3	3:F:623:PHE:HA	1.88	0.55
1:A:57:GLU:OE1	3:E:630:ARG:NH1	2.40	0.55
1:A:162:THR:OG1	1:A:205:ASN:HB2	2.07	0.55
3:E:605:LEU:HB3	3:E:607:PHE:CE2	2.42	0.55
3:E:580:THR:O	3:E:581:ASP:OD1	2.25	0.55
2:D:125:LEU:O	2:D:183:LYS:CE	2.55	0.55
2:D:1:ASP:N	2:D:95:PRO:HD2	2.22	0.55
2:D:198:HIS:CD2	2:D:200:THR:HG23	2.42	0.55
1:C:197:TRP:CZ2	1:C:219:ILE:HG22	2.42	0.55
3:E:536:LEU:HD22	3:E:637:ARG:NE	2.22	0.55
1:C:76:SER:HB2	1:C:78:THR:HG22	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:197:TRP:CA	1:C:199:SER:H	2.20	0.55
3:F:636:VAL:HG22	3:F:637:ARG:HG3	1.89	0.54
3:F:592:GLY:H	3:F:623:PHE:CB	2.20	0.54
1:C:150:LEU:C	1:C:150:LEU:HD23	2.28	0.54
2:B:189:HIS:CB	2:B:211:ARG:HD2	2.37	0.54
2:B:187:GLU:HA	2:B:211:ARG:NH2	2.09	0.54
2:D:193:THR:CG2	2:D:208:SER:HB3	2.31	0.54
1:C:102:ARG:HG3	1:C:102:ARG:NH1	2.23	0.54
2:B:119:PRO:HA	2:B:209:PHE:CZ	2.43	0.54
3:F:629:THR:HB	3:F:630:ARG:HH21	1.72	0.54
3:F:625:VAL:HG22	3:F:637:ARG:O	2.08	0.54
1:C:179:LEU:HD12	1:C:183:LEU:O	2.08	0.54
2:D:4:LEU:HD11	2:D:90:GLN:CB	2.38	0.54
1:A:47:TRP:CH2	1:A:49:GLY:HA2	2.43	0.54
2:B:124:GLN:NE2	2:B:131:SER:OG	2.40	0.54
1:C:199:SER:HB3	1:C:200:GLN:NE2	2.23	0.54
1:A:202:ILE:HG13	1:A:218:LYS:CE	2.38	0.54
3:E:514:ILE:HG22	3:F:512:HIS:O	2.08	0.53
1:A:161:LEU:HA	1:A:205:ASN:O	2.08	0.53
1:A:144:SER:HA	1:A:192:VAL:O	2.08	0.53
1:C:219:ILE:CD1	1:C:220:GLU:H	2.22	0.53
2:B:51:ALA:O	2:B:52:SER:HB3	2.07	0.53
3:F:614:LEU:HD12	3:F:637:ARG:NH2	2.24	0.53
1:A:205:ASN:HD22	1:A:215:VAL:HG12	1.74	0.53
1:C:203:THR:CG2	1:C:217:LYS:HE2	2.38	0.53
1:C:110:ASP:HB3	1:C:111:TYR:CD2	2.44	0.53
1:A:155:PHE:CD1	1:A:156:PRO:HA	2.44	0.53
1:C:110:ASP:HB3	1:C:111:TYR:HD2	1.74	0.53
3:E:524:LYS:HG2	3:E:525:ARG:H	1.73	0.52
2:B:115:VAL:HG21	2:B:205:ILE:HG12	1.92	0.52
3:E:606:THR:OG1	3:E:647:ASP:HA	2.09	0.52
2:B:180:THR:C	2:B:181:LEU:HD12	2.29	0.52
1:A:53:PRO:O	1:A:74:LYS:HE2	2.08	0.52
3:F:621:CYS:O	3:F:640:TYR:HB3	2.09	0.52
2:B:151:ASP:OD2	2:B:190:ASN:HB2	2.10	0.52
1:C:133:LEU:HB2	1:C:148:GLY:C	2.30	0.52
1:C:55:ASP:CG	1:C:57:GLU:HG2	2.29	0.52
1:C:170:SER:H	1:C:172:VAL:HG12	1.75	0.52
3:E:605:LEU:HB2	3:E:651:LEU:HB3	1.91	0.52
2:D:133:VAL:HG22	2:D:178:THR:HG23	1.92	0.52
3:F:553:ILE:HD11	3:F:649:ILE:HG21	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:136:VAL:HG11	2:D:119:PRO:CG	2.36	0.52
1:C:208:HIS:ND1	1:C:210:ALA:HB3	2.24	0.52
2:B:94:TRP:HB2	3:E:525:ARG:HG3	1.92	0.52
3:E:616:PRO:HB3	3:E:640:TYR:CZ	2.44	0.52
1:A:91:SER:O	1:A:92:ALA:HB2	2.09	0.52
1:A:175:PHE:CE2	2:B:176:SER:HB3	2.45	0.52
2:B:179:LEU:CD2	2:B:181:LEU:HD13	2.39	0.52
1:C:50:ARG:HD3	3:F:528:TYR:HB3	1.91	0.52
1:C:7:SER:HB3	1:C:21:SER:HB3	1.90	0.52
1:C:154:TYR:O	1:C:155:PHE:HB2	2.10	0.52
1:C:197:TRP:HA	1:C:199:SER:N	2.25	0.51
1:C:150:LEU:HD21	1:C:152:LYS:HB2	1.92	0.51
2:B:155:ARG:HH21	2:B:181:LEU:HD11	1.76	0.51
2:B:90:GLN:NE2	2:B:93:SER:N	2.58	0.51
3:E:523:MET:SD	3:E:533:ALA:HA	2.50	0.51
3:F:570:LEU:CD2	3:F:577:ARG:HE	2.24	0.51
3:E:558:PHE:CZ	3:E:651:LEU:HD13	2.44	0.51
1:C:133:LEU:HD11	1:C:150:LEU:HB2	1.93	0.51
1:C:171:GLY:O	1:C:190:VAL:HA	2.11	0.51
2:B:94:TRP:CG	3:E:525:ARG:HD2	2.46	0.51
1:C:18:VAL:HG22	1:C:83:LEU:HB2	1.91	0.51
2:B:1:ASP:N	3:E:571:TRP:HZ2	2.08	0.51
1:A:164:ASN:OD1	1:A:202:ILE:HA	2.11	0.51
3:F:602:ASN:O	3:F:603:ASN:C	2.49	0.51
3:F:599:ASP:O	3:F:600:LYS:CB	2.58	0.51
3:E:648:ASN:N	3:E:648:ASN:HD22	2.07	0.51
3:F:616:PRO:HB2	3:F:640:TYR:CD1	2.45	0.51
2:B:189:HIS:CE1	2:B:211:ARG:HG3	2.45	0.51
2:B:118:PHE:O	2:B:132:VAL:HG23	2.11	0.51
1:C:2:VAL:HG22	1:C:27:TYR:HB3	1.92	0.51
2:B:130:ALA:HB3	2:B:181:LEU:O	2.10	0.51
1:C:52:TYR:HD1	1:C:55:ASP:HB3	1.75	0.51
1:C:55:ASP:OD2	1:C:57:GLU:CG	2.59	0.51
3:F:628:ARG:NH1	3:F:633:GLU:CD	2.64	0.51
2:B:167:ASP:HB3	2:B:170:ASP:OD1	2.11	0.51
2:B:127:SER:C	2:B:129:GLY:H	2.14	0.51
1:A:160:THR:HG23	1:A:207:ALA:HB3	1.92	0.51
1:C:219:ILE:CG1	1:C:220:GLU:H	2.24	0.51
1:A:209:PRO:C	1:A:211:SER:N	2.64	0.51
3:E:579:CYS:SG	3:E:583:LEU:O	2.69	0.51
1:C:168:LEU:HG	1:C:168:LEU:O	2.10	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:193:THR:HG22	2:B:208:SER:CB	2.39	0.50
3:E:536:LEU:HD22	3:E:637:ARG:HE	1.76	0.50
1:C:20:ILE:HG22	1:C:21:SER:N	2.26	0.50
2:B:1:ASP:N	3:E:571:TRP:CZ2	2.79	0.50
3:F:624:ASP:HB3	3:F:635:VAL:CG1	2.36	0.50
1:C:102:ARG:O	1:C:103:TYR:CB	2.59	0.50
1:A:189:SER:HB2	2:B:135:PHE:CD2	2.47	0.50
2:D:183:LYS:O	2:D:187:GLU:HG2	2.11	0.50
2:B:159:VAL:CG1	2:B:179:LEU:HG	2.42	0.50
3:E:553:ILE:HD12	3:E:558:PHE:CD2	2.47	0.50
1:C:108:ALA:HB2	2:D:34:HIS:CE1	2.47	0.50
1:A:146:THR:HG22	1:A:191:THR:HG23	1.93	0.50
1:C:12:VAL:HG11	1:C:18:VAL:CG1	2.42	0.50
3:E:526:SER:OG	3:E:530:GLN:HB3	2.11	0.50
1:C:60:TYR:OH	1:C:69:THR:HA	2.11	0.50
2:D:3:LEU:HD22	2:D:5:THR:HG23	1.93	0.50
1:A:189:SER:HB2	2:B:135:PHE:CE2	2.47	0.50
1:C:162:THR:OG1	1:C:205:ASN:HB2	2.12	0.50
3:F:616:PRO:CG	3:F:640:TYR:CG	2.90	0.50
1:C:128:PRO:HA	1:C:154:TYR:HB3	1.93	0.50
1:A:181:SER:OG	1:A:182:ASP:N	2.40	0.49
2:B:149:LYS:C	2:B:150:ILE:HG13	2.32	0.49
1:C:154:TYR:HE1	1:C:157:GLU:HG3	1.76	0.49
3:F:569:SER:HB2	3:F:575:PHE:CD1	2.47	0.49
1:C:103:TYR:HA	3:F:530:GLN:HE21	1.77	0.49
1:C:103:TYR:HA	3:F:530:GLN:NE2	2.27	0.49
2:B:196:ALA:O	2:B:204:PRO:HA	2.12	0.49
1:A:33:TRP:CE3	1:A:50:ARG:HD2	2.47	0.49
3:F:568:SER:HB3	3:F:577:ARG:HB3	1.93	0.49
1:C:62:GLU:HG2	1:C:65:LYS:HD2	1.94	0.49
2:B:140:TYR:CG	2:B:141:PRO:HA	2.47	0.49
1:C:203:THR:HG22	1:C:217:LYS:HE2	1.94	0.49
3:F:603:ASN:O	3:F:604:TYR:CD1	2.65	0.49
3:F:612:LEU:HD22	3:F:639:LEU:CD1	2.41	0.49
1:C:106:TYR:OH	3:F:532:ILE:HD13	2.12	0.49
3:F:536:LEU:N	3:F:536:LEU:HD13	2.25	0.49
1:C:87:THR:HG22	1:C:89:VAL:H	1.78	0.49
1:A:144:SER:HB3	1:A:191:THR:CG2	2.42	0.49
3:E:553:ILE:HG13	3:E:553:ILE:O	2.12	0.49
1:A:33:TRP:CZ3	1:A:52:TYR:HB2	2.48	0.49
2:B:135:PHE:C	2:B:136:LEU:HD12	2.33	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:51:ILE:HD12	1:C:70:LEU:O	2.13	0.48
2:D:190:ASN:ND2	2:D:211:ARG:HG3	2.27	0.48
2:B:120:PRO:HG2	2:B:186:TYR:CE1	2.47	0.48
1:A:73:ASP:HB3	1:A:76:SER:HB3	1.96	0.48
2:B:89:GLN:HB2	2:B:98:PHE:CE1	2.48	0.48
2:B:209:PHE:HD1	2:B:210:ASN:O	1.96	0.48
1:C:169:SER:O	1:C:170:SER:CB	2.61	0.48
3:F:548:THR:O	3:F:549:ASP:CB	2.61	0.48
1:A:163:TRP:CZ3	1:A:204:CYS:HB3	2.48	0.48
2:D:193:THR:HG22	2:D:208:SER:CB	2.38	0.48
1:C:52:TYR:CD1	1:C:55:ASP:HB3	2.48	0.48
2:B:184:ASP:C	2:B:186:TYR:H	2.15	0.48
2:B:183:LYS:O	2:B:187:GLU:HG2	2.13	0.48
2:B:146:VAL:CG1	2:B:147:LYS:N	2.76	0.48
3:F:628:ARG:NH1	3:F:628:ARG:HG2	2.26	0.48
3:E:622:LYS:HE3	3:E:638:SER:HB3	1.96	0.48
3:F:596:PHE:HE2	3:F:643:TYR:CZ	2.32	0.48
2:B:115:VAL:HA	2:B:135:PHE:O	2.14	0.47
1:C:175:PHE:CE2	2:D:176:SER:HB3	2.49	0.47
1:C:50:ARG:HG2	1:C:59:ASN:HB2	1.96	0.47
3:F:570:LEU:HD21	3:F:577:ARG:NH2	2.24	0.47
1:A:45:LEU:HD12	1:A:45:LEU:N	2.29	0.47
2:B:150:ILE:HD13	2:B:192:TYR:HE1	1.79	0.47
3:E:574:ILE:O	3:E:576:LYS:N	2.48	0.47
2:B:10:ILE:HD13	2:B:142:LYS:HZ3	1.79	0.47
1:C:133:LEU:HB2	1:C:148:GLY:CA	2.45	0.47
2:D:29:ILE:HA	2:D:92:ASP:OD2	2.15	0.47
1:A:18:VAL:HG12	1:A:83:LEU:HB2	1.97	0.47
1:C:125:THR:HA	1:C:156:PRO:HD2	1.93	0.47
1:A:161:LEU:HD21	1:A:188:SER:HB2	1.97	0.47
1:C:50:ARG:NH1	3:F:528:TYR:HB2	2.30	0.47
1:C:210:ALA:O	1:C:212:SER:N	2.48	0.47
3:E:577:ARG:CG	3:E:578:ASN:N	2.78	0.47
1:A:161:LEU:HD12	1:A:205:ASN:O	2.15	0.47
1:A:124:LYS:O	1:A:126:THR:HG23	2.14	0.47
1:C:130:VAL:HG12	1:C:214:LYS:HD3	1.96	0.47
1:C:193:THR:HB	1:C:196:THR:OG1	2.15	0.47
1:A:164:ASN:HD21	1:A:202:ILE:HG22	1.78	0.47
2:B:126:THR:C	2:B:128:GLY:H	2.18	0.47
2:B:128:GLY:HA2	2:B:183:LYS:HE2	1.96	0.47
1:A:157:GLU:CG	1:A:158:PRO:HA	2.45	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:513:THR:HG22	3:E:555:SER:CB	2.44	0.47
1:C:61:SER:HB3	1:C:64:PHE:HD1	1.80	0.47
3:E:539:ILE:CG1	3:E:614:LEU:HD11	2.45	0.47
2:B:94:TRP:CB	3:E:525:ARG:HG3	2.45	0.47
3:F:631:THR:HG23	3:F:631:THR:O	2.15	0.47
3:F:546:ASN:O	3:F:546:ASN:CG	2.54	0.47
3:F:617:VAL:HG22	3:F:618:GLY:N	2.29	0.47
2:D:17:GLU:O	2:D:78:VAL:HG23	2.14	0.47
2:B:2:ILE:HG13	2:B:26:SER:HB2	1.96	0.46
3:E:605:LEU:HB3	3:E:607:PHE:HE2	1.80	0.46
3:E:619:ALA:HA	3:E:642:ILE:HG12	1.97	0.46
3:E:576:LYS:O	3:E:577:ARG:HB2	2.14	0.46
1:A:55:ASP:OD2	1:A:57:GLU:HG2	2.15	0.46
1:C:36:TRP:CD1	1:C:70:LEU:HD22	2.50	0.46
3:F:596:PHE:CE2	3:F:643:TYR:CE1	3.03	0.46
3:E:653:PRO:CA	3:E:654:ARG:CB	2.94	0.46
2:B:125:LEU:HD12	2:B:129:GLY:O	2.15	0.46
2:B:151:ASP:N	2:B:190:ASN:OD1	2.49	0.46
2:B:196:ALA:HB3	2:B:205:ILE:CG2	2.46	0.46
1:C:144:SER:HB2	1:C:191:THR:CG2	2.45	0.46
2:D:157:ASN:N	2:D:157:ASN:OD1	2.49	0.46
2:D:2:ILE:HD11	2:D:27:GLN:HE21	1.81	0.46
1:C:87:THR:HG22	1:C:88:SER:N	2.31	0.46
3:E:621:CYS:SG	3:E:641:VAL:HG13	2.56	0.46
2:D:18:ARG:NH1	2:D:76:ASN:HA	2.31	0.46
3:F:582:VAL:CG1	3:F:582:VAL:O	2.64	0.46
1:C:192:VAL:HG11	1:C:202:ILE:CD1	2.44	0.46
2:B:135:PHE:CD1	2:B:135:PHE:N	2.84	0.46
1:A:67:LYS:HE3	1:A:90:ASP:CG	2.36	0.46
3:E:574:ILE:C	3:E:576:LYS:H	2.20	0.45
1:C:186:LEU:HD12	1:C:186:LEU:C	2.36	0.45
1:C:68:ALA:HA	1:C:82:HIS:O	2.16	0.45
1:C:155:PHE:CE1	1:C:156:PRO:HA	2.51	0.45
2:D:119:PRO:HG3	2:D:209:PHE:CG	2.51	0.45
3:E:525:ARG:NH2	3:E:581:ASP:OD1	2.49	0.45
2:B:162:SER:O	2:B:175:MET:HG3	2.17	0.45
3:E:513:THR:HG22	3:E:555:SER:HB3	1.99	0.45
2:D:13:VAL:HG21	2:D:78:VAL:HG21	1.98	0.45
2:D:2:ILE:HG23	2:D:26:SER:HB2	1.97	0.45
1:A:67:LYS:HE3	1:A:90:ASP:OD2	2.16	0.45
1:C:30:SER:HA	1:C:53:PRO:CB	2.32	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:193:THR:HA	2:D:208:SER:HB3	1.98	0.45
1:C:157:GLU:OE1	1:C:184:TYR:CE1	2.69	0.45
3:F:568:SER:OG	3:F:583:LEU:HD23	2.16	0.45
2:B:151:ASP:OD1	2:B:190:ASN:HB2	2.12	0.45
1:A:161:LEU:CD2	1:A:188:SER:HB2	2.46	0.45
2:D:107:LYS:HB2	2:D:107:LYS:HE3	1.67	0.45
1:C:128:PRO:HG3	1:C:208:HIS:CD2	2.51	0.45
1:A:67:LYS:HE3	1:A:90:ASP:OD1	2.17	0.45
2:D:115:VAL:HA	2:D:135:PHE:O	2.16	0.45
2:D:117:ILE:HD13	2:D:192:TYR:O	2.17	0.45
1:C:136:VAL:C	1:C:138:GLY:CA	2.85	0.44
1:C:133:LEU:HD11	1:C:150:LEU:CB	2.47	0.44
1:A:11:LEU:HD11	1:A:156:PRO:HB3	1.98	0.44
2:D:120:PRO:HD3	2:D:132:VAL:HG22	1.98	0.44
3:F:584:ASP:HB3	3:F:630:ARG:HH22	1.83	0.44
1:A:40:ARG:HG3	1:A:41:PRO:CD	2.42	0.44
1:C:176:PRO:HG2	2:D:163:TRP:O	2.16	0.44
3:F:608:ASN:HA	3:F:648:ASN:HB2	1.98	0.44
1:C:47:TRP:CH2	1:C:49:GLY:HA2	2.53	0.44
1:A:175:PHE:CD2	2:B:176:SER:HB3	2.52	0.44
3:F:596:PHE:HE2	3:F:643:TYR:CE1	2.35	0.44
3:F:613:SER:HB3	3:F:616:PRO:HG3	2.00	0.44
3:F:622:LYS:HE2	3:F:638:SER:OG	2.18	0.44
1:C:179:LEU:HD13	1:C:184:TYR:CE2	2.52	0.44
1:A:147:LEU:HB2	1:A:218:LYS:NZ	2.31	0.44
3:E:512:HIS:C	3:E:512:HIS:CD2	2.90	0.44
3:F:602:ASN:CG	3:F:603:ASN:H	2.20	0.44
3:F:653:PRO:HB2	3:F:654:ARG:H	1.36	0.44
1:A:67:LYS:CE	1:A:90:ASP:OD2	2.66	0.44
3:E:520:ASP:N	3:E:538:ASN:O	2.50	0.44
3:F:579:CYS:HB3	3:F:583:LEU:O	2.17	0.44
2:D:36:TYR:HA	2:D:45:ARG:O	2.17	0.44
1:C:135:PRO:HD3	1:C:147:LEU:HD13	2.00	0.44
2:D:38:GLN:O	2:D:84:ALA:HB1	2.18	0.44
2:B:115:VAL:HG23	2:B:207:LYS:CG	2.47	0.44
2:B:150:ILE:HD13	2:B:192:TYR:CE1	2.52	0.43
2:D:33:ILE:CG2	2:D:51:ALA:HB2	2.44	0.43
2:D:179:LEU:HG	2:D:181:LEU:HD13	1.99	0.43
3:F:561:TYR:CZ	5:F:701:NAG:H82	2.54	0.43
3:E:639:LEU:HD23	3:E:639:LEU:HA	1.80	0.43
2:B:39:ARG:NH2	4:B:301:ACY:OXT	2.51	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:72:THR:HG22	2:D:74:ASN:ND2	2.32	0.43
3:E:513:THR:O	3:E:558:PHE:HA	2.18	0.43
3:F:592:GLY:CA	3:F:623:PHE:HA	2.47	0.43
3:E:514:ILE:HG23	3:F:512:HIS:HB3	2.00	0.43
3:E:625:VAL:HG23	3:E:625:VAL:O	2.19	0.43
2:B:32:SER:O	2:B:91:THR:HG22	2.19	0.43
3:F:534:SER:O	3:F:636:VAL:HG23	2.18	0.43
2:B:179:LEU:CD2	2:B:181:LEU:CD1	2.97	0.43
2:D:198:HIS:HD2	2:D:200:THR:HG23	1.80	0.43
2:D:170:ASP:OD1	2:D:172:THR:HG23	2.19	0.43
1:C:39:GLN:HB2	1:C:45:LEU:HD23	2.01	0.43
3:F:612:LEU:HD23	3:F:612:LEU:HA	1.85	0.43
3:E:570:LEU:HB3	3:E:571:TRP:H	1.46	0.43
2:D:15:PRO:HD3	2:D:106:LEU:HD11	2.01	0.43
1:C:202:ILE:O	1:C:218:LYS:HE3	2.18	0.43
1:A:197:TRP:CD2	1:A:198:PRO:HA	2.54	0.43
3:F:600:LYS:HE3	3:F:653:PRO:HB2	2.01	0.43
3:F:567:LYS:HD3	3:F:567:LYS:HA	1.89	0.43
2:B:181:LEU:HG	2:B:185:GLU:CD	2.39	0.43
3:F:570:LEU:HA	3:F:570:LEU:HD12	1.71	0.43
1:A:60:TYR:O	2:B:94:TRP:HH2	2.02	0.43
3:F:517:ILE:HD11	3:F:560:VAL:HG11	2.01	0.43
1:A:133:LEU:HB2	1:A:148:GLY:O	2.19	0.43
1:A:150:LEU:HD23	1:A:150:LEU:C	2.39	0.43
1:A:131:TYR:HA	1:A:132:PRO:HD3	1.84	0.43
3:F:623:PHE:CE2	3:F:639:LEU:HD12	2.54	0.42
1:A:196:THR:HG23	1:A:200:GLN:HB2	2.00	0.42
1:A:150:LEU:HD21	1:A:152:LYS:HB2	2.01	0.42
2:B:106:LEU:HA	2:B:106:LEU:HD23	1.73	0.42
2:D:6:GLN:HA	2:D:22:SER:O	2.19	0.42
1:C:153:GLY:HA2	1:C:183:LEU:HB3	2.00	0.42
3:E:597:SER:O	3:E:598:PHE:CB	2.67	0.42
1:A:146:THR:HG22	1:A:191:THR:CG2	2.49	0.42
1:A:219:ILE:O	1:A:219:ILE:HG22	2.20	0.42
3:E:598:PHE:O	3:E:598:PHE:HD2	2.01	0.42
3:E:550:VAL:HA	3:E:612:LEU:O	2.19	0.42
3:F:574:ILE:HG23	3:F:574:ILE:O	2.19	0.42
2:B:21:LEU:HB2	2:B:73:LEU:HB3	2.01	0.42
3:E:553:ILE:HD11	3:E:649:ILE:HG21	2.01	0.42
1:A:51:ILE:O	1:A:53:PRO:HD3	2.20	0.42
1:A:128:PRO:HA	1:A:154:TYR:HB3	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:582:VAL:C	3:F:584:ASP:H	2.23	0.42
3:F:581:ASP:HB2	3:F:582:VAL:H	1.54	0.42
3:F:582:VAL:HG12	3:F:582:VAL:O	2.20	0.42
3:F:628:ARG:CG	3:F:628:ARG:HH11	2.25	0.42
1:A:45:LEU:CD1	1:A:45:LEU:N	2.82	0.42
1:A:38:LYS:HB2	1:A:48:ILE:HD11	2.02	0.42
2:D:190:ASN:CB	2:D:211:ARG:HD3	2.42	0.42
1:A:105:PRO:HB3	3:E:530:GLN:HG2	2.01	0.42
3:F:592:GLY:N	3:F:623:PHE:HA	2.34	0.42
2:B:159:VAL:HG12	2:B:179:LEU:HG	2.01	0.42
2:B:90:GLN:NE2	2:B:92:ASP:N	2.65	0.42
3:E:536:LEU:HA	3:E:536:LEU:HD22	1.72	0.42
1:A:159:VAL:CG1	1:A:186:LEU:HD21	2.50	0.42
3:F:566:CYS:SG	3:F:585:ALA:HB3	2.60	0.42
3:E:591:THR:HG23	3:E:591:THR:O	2.19	0.42
2:B:135:PHE:HB3	2:B:137:ASN:ND2	2.35	0.42
3:E:598:PHE:O	3:E:598:PHE:CD2	2.72	0.42
3:E:612:LEU:N	3:E:612:LEU:CD1	2.83	0.42
3:E:596:PHE:HB3	3:E:605:LEU:HD11	2.01	0.42
3:F:651:LEU:HG	3:F:652:VAL:N	2.34	0.42
1:A:101:TYR:O	1:A:102:ARG:C	2.58	0.42
3:F:518:THR:HG23	3:F:540:THR:OG1	2.20	0.42
3:F:613:SER:O	3:F:616:PRO:HG3	2.20	0.42
2:D:190:ASN:CG	2:D:190:ASN:O	2.58	0.42
2:D:131:SER:HA	2:D:179:LEU:O	2.19	0.42
3:E:593:THR:OG1	3:E:621:CYS:HB2	2.20	0.42
3:E:594:CYS:HA	3:E:595:PRO:HD3	1.84	0.42
2:D:149:LYS:HB3	2:D:154:GLU:HA	2.02	0.42
3:E:537:SER:HB2	3:E:636:VAL:HG11	2.01	0.42
3:F:580:THR:CG2	3:F:581:ASP:H	2.22	0.42
3:E:614:LEU:HD12	3:E:614:LEU:HA	1.87	0.42
2:D:189:HIS:O	2:D:190:ASN:CB	2.68	0.42
3:F:553:ILE:CD1	3:F:649:ILE:HG21	2.49	0.42
2:B:135:PHE:HD1	2:B:135:PHE:N	2.18	0.41
1:C:62:GLU:HG2	1:C:65:LYS:CD	2.49	0.41
1:C:143:SER:HA	1:C:144:SER:HA	1.61	0.41
3:F:594:CYS:HA	3:F:595:PRO:HD3	1.79	0.41
1:C:197:TRP:CH2	1:C:219:ILE:HA	2.56	0.41
2:B:45:ARG:HA	2:B:46:PRO:HD3	1.77	0.41
1:C:161:LEU:C	1:C:161:LEU:HD23	2.40	0.41
2:D:190:ASN:HD21	2:D:210:ASN:C	2.24	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:146:THR:HG22	1:A:191:THR:OG1	2.21	0.41
3:E:645:GLU:O	3:E:645:GLU:HG3	2.21	0.41
1:C:102:ARG:HG2	1:C:103:TYR:CE2	2.56	0.41
2:B:146:VAL:HG12	2:B:147:LYS:N	2.34	0.41
1:A:147:LEU:HD12	1:A:202:ILE:HD11	2.01	0.41
1:A:147:LEU:O	1:A:190:VAL:HG12	2.19	0.41
3:F:606:THR:HB	3:F:647:ASP:HA	2.02	0.41
3:F:583:LEU:HD12	3:F:583:LEU:HA	1.82	0.41
1:A:201:SER:HB3	1:A:217:LYS:HE3	2.02	0.41
2:D:130:ALA:O	2:D:181:LEU:N	2.45	0.41
2:D:179:LEU:HG	2:D:181:LEU:CD1	2.50	0.41
1:A:180:GLN:HG2	1:A:183:LEU:O	2.21	0.41
1:A:110:ASP:HB3	1:A:111:TYR:HD2	1.84	0.41
3:F:638:SER:CB	3:F:640:TYR:CE2	3.03	0.41
2:B:189:HIS:CG	2:B:190:ASN:N	2.88	0.41
1:C:6:GLN:HA	1:C:21:SER:O	2.20	0.41
3:F:617:VAL:CG2	3:F:618:GLY:N	2.83	0.41
2:B:189:HIS:NE2	2:B:210:ASN:CB	2.71	0.41
3:E:608:ASN:CG	3:E:646:GLY:HA3	2.41	0.41
1:C:218:LYS:HB3	1:C:219:ILE:H	1.41	0.41
1:A:145:VAL:O	1:A:191:THR:HA	2.21	0.41
1:A:159:VAL:HG23	1:A:207:ALA:O	2.21	0.41
3:E:622:LYS:HE3	3:E:638:SER:CB	2.51	0.41
3:F:539:ILE:HG23	3:F:614:LEU:CD1	2.43	0.41
3:F:536:LEU:CD2	3:F:575:PHE:HE2	2.34	0.41
1:C:55:ASP:OD1	1:C:55:ASP:C	2.59	0.41
2:D:1:ASP:H1	2:D:95:PRO:CD	2.34	0.41
2:D:1:ASP:H2	2:D:95:PRO:HD2	1.86	0.41
1:A:104:ASP:HA	1:A:105:PRO:HD3	1.79	0.41
1:A:89:VAL:O	1:A:89:VAL:HG12	2.21	0.41
2:D:203:SER:HA	2:D:204:PRO:HD3	1.94	0.41
3:F:534:SER:OG	3:F:536:LEU:HD22	2.21	0.41
1:C:197:TRP:CG	1:C:198:PRO:HA	2.56	0.41
1:A:176:PRO:HG2	2:B:163:TRP:O	2.21	0.41
3:F:601:LEU:HD12	3:F:601:LEU:HA	1.80	0.41
3:F:539:ILE:O	3:F:614:LEU:HD11	2.21	0.40
2:D:124:GLN:HG2	2:D:130:ALA:HA	2.01	0.40
1:C:145:VAL:O	1:C:191:THR:HA	2.22	0.40
1:A:36:TRP:O	1:A:48:ILE:HB	2.21	0.40
3:F:536:LEU:HD21	3:F:575:PHE:HE2	1.86	0.40
3:F:596:PHE:CE2	3:F:643:TYR:CZ	3.09	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:178:VAL:HG12	1:A:185:THR:O	2.21	0.40
1:A:35:ASN:HB3	1:A:47:TRP:HE1	1.85	0.40
2:D:40:THR:C	2:D:42:GLY:H	2.25	0.40
3:F:623:PHE:CZ	3:F:641:VAL:HG11	2.56	0.40
3:E:616:PRO:HB3	3:E:640:TYR:CE1	2.55	0.40
3:E:530:GLN:HA	3:E:531:PRO:HD3	1.84	0.40
2:D:82:ASP:HA	7:D:401:HOH:O	2.21	0.40
2:B:136:LEU:N	2:B:136:LEU:HD12	2.37	0.40
3:F:560:VAL:HG12	3:F:561:TYR:N	2.37	0.40
1:C:65:LYS:C	1:C:67:LYS:H	2.24	0.40
2:B:123:GLU:N	2:B:123:GLU:OE1	2.54	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	210/221 (95%)	186 (89%)	17 (8%)	7 (3%)	5	26
1	C	218/221 (99%)	186 (85%)	18 (8%)	14 (6%)	2	9
2	B	211/214 (99%)	177 (84%)	22 (10%)	12 (6%)	2	12
2	D	209/214 (98%)	193 (92%)	14 (7%)	2 (1%)	19	61
3	E	146/209 (70%)	121 (83%)	18 (12%)	7 (5%)	3	17
3	F	146/209 (70%)	119 (82%)	24 (16%)	3 (2%)	9	40
All	All	1140/1288 (88%)	982 (86%)	113 (10%)	45 (4%)	4	21

All (45) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	52	SER

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Mol	Chain	Res	Type
2	B	131	SER
2	B	153	SER
2	B	154	GLU
2	B	212	ASN
1	C	137	CYS
1	C	170	SER
1	C	181	SER
1	C	211	SER
1	C	219	ILE
3	E	549	ASP
3	F	603	ASN
3	F	653	PRO
1	A	102	ARG
2	B	51	ALA
2	B	156	GLN
2	B	158	GLY
1	C	142	GLY
2	D	52	SER
2	D	190	ASN
3	E	577	ARG
3	E	602	ASN
3	F	600	LYS
1	A	144	SER
1	A	210	ALA
2	B	127	SER
1	C	158	PRO
3	E	584	ASP
2	B	151	ASP
1	C	138	GLY
1	C	155	PHE
3	E	575	PHE
1	A	168	LEU
1	A	181	SER
2	B	190	ASN
1	C	103	TYR
1	C	165	SER
1	C	194	SER
3	E	544	GLN
3	E	545	ASP
2	B	143	ASP
1	C	195	SER
1	C	210	ALA

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Mol	Chain	Res	Type
1	A	209	PRO
1	A	105	PRO

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	181/187 (97%)	174 (96%)	7 (4%)	39	77
1	C	186/187 (100%)	173 (93%)	13 (7%)	19	55
2	B	190/191 (100%)	183 (96%)	7 (4%)	41	79
2	D	188/191 (98%)	181 (96%)	7 (4%)	41	79
3	E	135/187 (72%)	122 (90%)	13 (10%)	10	38
3	F	135/187 (72%)	121 (90%)	14 (10%)	9	32
All	All	1015/1130 (90%)	954 (94%)	61 (6%)	24	62

All (61) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	40	ARG
1	A	55	ASP
1	A	110	ASP
1	A	143	SER
1	A	188	SER
1	A	206	VAL
1	A	213	THR
2	B	1	ASP
2	B	7	SER
2	B	153	SER
2	B	154	GLU
2	B	159	VAL
2	B	191	SER
2	B	202	THR
1	C	21	SER
1	C	62	GLU

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Mol	Chain	Res	Type
1	C	93	VAL
1	C	102	ARG
1	C	110	ASP
1	C	139	ASP
1	C	170	SER
1	C	193	THR
1	C	195	SER
1	C	196	THR
1	C	211	SER
1	C	218	LYS
1	C	219	ILE
2	D	3	LEU
2	D	43	SER
2	D	124	GLN
2	D	127	SER
2	D	155	ARG
2	D	157	ASN
2	D	159	VAL
3	E	521	LEU
3	E	535	THR
3	E	536	LEU
3	E	550	VAL
3	E	572	ASP
3	E	581	ASP
3	E	584	ASP
3	E	594	CYS
3	E	606	THR
3	E	632	ASN
3	E	641	VAL
3	E	647	ASP
3	E	648	ASN
3	F	517	ILE
3	F	518	THR
3	F	525	ARG
3	F	536	LEU
3	F	539	ILE
3	F	547	ASN
3	F	570	LEU
3	F	581	ASP
3	F	603	ASN
3	F	614	LEU
3	F	617	VAL

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Mol	Chain	Res	Type
3	F	623	PHE
3	F	630	ARG
3	F	640	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	35	ASN
1	A	205	ASN
2	B	90	GLN
2	B	137	ASN
2	B	210	ASN
1	C	59	ASN
1	C	200	GLN
2	D	27	GLN
2	D	37	GLN
2	D	198	HIS
2	D	210	ASN
3	E	648	ASN
3	F	557	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

8 carbohydrates are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	NAG	E	701	3,5	14,14,15	0.78	0	15,19,21	1.17	2 (13%)
5	NAG	E	702	5	14,14,15	0.61	0	15,19,21	0.68	0
5	NAG	E	703	3,5	14,14,15	0.61	0	15,19,21	1.86	3 (20%)
5	NAG	E	704	5	14,14,15	0.55	0	15,19,21	1.43	1 (6%)
5	NAG	F	701	3,5	14,14,15	0.79	0	15,19,21	1.05	1 (6%)
5	NAG	F	702	5	14,14,15	0.62	0	15,19,21	0.63	0
5	NAG	F	703	3,5	14,14,15	0.67	0	15,19,21	1.06	1 (6%)
5	NAG	F	704	5	14,14,15	0.45	0	15,19,21	1.09	2 (13%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	E	701	3,5	-	0/6/23/26	0/1/1/1
5	NAG	E	702	5	-	0/6/23/26	0/1/1/1
5	NAG	E	703	3,5	-	0/6/23/26	0/1/1/1
5	NAG	E	704	5	-	0/6/23/26	0/1/1/1
5	NAG	F	701	3,5	-	0/6/23/26	0/1/1/1
5	NAG	F	702	5	-	0/6/23/26	0/1/1/1
5	NAG	F	703	3,5	-	0/6/23/26	0/1/1/1
5	NAG	F	704	5	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	F	701	NAG	C2-N2-C7	-2.39	119.97	123.04
5	E	703	NAG	C6-C5-C4	-2.05	107.96	113.02
5	F	704	NAG	C2-N2-C7	-2.02	120.44	123.04
5	E	701	NAG	O4-C4-C3	2.09	115.05	110.34
5	F	703	NAG	C4-C3-C2	2.26	114.74	111.23
5	F	704	NAG	C1-O5-C5	2.69	115.66	112.25
5	E	701	NAG	C1-O5-C5	2.80	115.80	112.25
5	E	703	NAG	C3-C4-C5	4.30	117.70	110.20
5	E	704	NAG	C1-O5-C5	4.63	118.13	112.25
5	E	703	NAG	C1-O5-C5	4.75	118.28	112.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	F	701	NAG	3	0
5	F	702	NAG	1	0

5.6 Ligand geometry

3 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	ACY	B	301	-	1,3,3	0.87	0	0,3,3	0.00	-
4	ACY	D	301	-	1,3,3	0.62	0	0,3,3	0.00	-
6	NAG	F	705	3	14,14,15	0.52	0	15,19,21	0.83	1 (6%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	ACY	B	301	-	-	0/0/0/0	0/0/0/0
4	ACY	D	301	-	-	0/0/0/0	0/0/0/0
6	NAG	F	705	3	1/1/5/7	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	F	705	NAG	C2-N2-C7	-2.24	120.16	123.04

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
6	F	705	NAG	C1

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	301	ACY	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	214/221 (96%)	-0.40	1 (0%) 91 76	34, 67, 121, 166	1 (0%)
1	C	220/221 (99%)	-0.27	3 (1%) 78 51	43, 83, 118, 157	3 (1%)
2	B	213/214 (99%)	-0.21	3 (1%) 78 51	39, 77, 158, 170	4 (1%)
2	D	211/214 (98%)	-0.50	1 (0%) 91 76	32, 62, 114, 127	0
3	E	148/209 (70%)	-0.23	3 (2%) 68 39	54, 83, 146, 185	0
3	F	148/209 (70%)	-0.30	0 100 100	50, 87, 136, 187	0
All	All	1154/1288 (89%)	-0.32	11 (0%) 84 60	32, 77, 138, 187	8 (0%)

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	212	ASN	4.3
2	B	134	CYS	3.7
1	C	137	CYS	3.7
3	E	571	TRP	3.1
3	E	546	ASN	3.1
1	C	167	SER	2.9
1	A	141	THR	2.9
3	E	602	ASN	2.8
2	D	190	ASN	2.6
2	B	157	ASN	2.4
1	C	213	THR	2.3

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
5	NAG	F	701	14/15	0.94	0.32	2.44	96,102,106,114	0
5	NAG	E	701	14/15	0.91	0.21	0.24	85,94,106,118	0
5	NAG	F	703	14/15	0.95	0.16	0.10	91,96,106,111	0
5	NAG	E	703	14/15	0.91	0.15	-0.32	74,116,123,134	0
5	NAG	E	702	14/15	0.89	0.33	-	130,137,143,145	0
5	NAG	F	704	14/15	0.90	0.33	-	117,119,121,122	0
5	NAG	E	704	14/15	0.91	0.30	-	137,147,152,152	0
5	NAG	F	702	14/15	0.80	0.43	-	123,126,134,135	0

6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
4	ACY	B	301	4/4	0.93	0.18	2.02	55,69,73,78	0
4	ACY	D	301	4/4	0.96	0.14	0.42	51,59,59,65	0
6	NAG	F	705	14/15	0.82	0.42	-	150,154,156,157	0

6.5 Other polymers ⓘ

There are no such residues in this entry.